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PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

NEWS 1 Web Page for STN Seminar Schedule - N. America

NEWS 2 AUG 10 Time limit for inactive STN sessions doubles to 40 minutes

NEWS 3 AUG 18 COMPENDEX indexing changed for the Corporate Source (CS) field

NEWS 4 AUG 24 ENCOMPLIT/ENCOMPLIT2 reloaded and enhanced

NEWS 5 AUG 24 CA/CAplus enhanced with legal status information for U.S. patents

NEWS 6 SEP 09 50 Millionth Unique Chemical Substance Recorded in CAS REGISTRY

NEWS 7 SEP 11 WPIDS, WPINDEX, and WPIX now include Japanese FTERM thesaurus

NEWS 8 OCT 21 Derwent World Patents Index Coverage of Indian and Taiwanese Content Expanded

NEWS 9 OCT 21 Derwent World Patents Index enhanced with human translated claims for Chinese Applications and Utility Models

NEWS 10 OCT 27 Free display of legal status information in CA/CAplus, USPATFULL, and USPAT2 in the month of November.

NEWS EXPRESS MAY 26 09 CURRENT WINDOWS VERSION IS V8.4, AND CURRENT DISCOVER FILE IS DATED 06 APRIL 2009.

NEWS HOURS STN Operating Hours Plus Help Desk Availability NEWS LOGIN Welcome Banner and News Items

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 14:23:13 ON 18 NOV 2009

=> file registry COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.22 0.22

FULL ESTIMATED COST

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8 DICTIONARY FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 26, 2009.

Please note that search-term pricing does apply when conducting ${\tt SmartSELECT}$ searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

=> file hcaplus COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.48 0.70

FULL ESTIMATED COST

FILE 'HCAPLUS' ENTERED AT 14:24:18 ON 18 NOV 2009 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2009 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 18 Nov 2009 VOL 151 ISS 21
FILE LAST UPDATED: 17 Nov 2009 (20091117/ED)
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Aug 2009
USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Aug 2009

HCAplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2009.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

During November, try the new LSUS format of legal status information in the ${\rm CA/CAplus}$ family databases for free! Complete details on the

number of free displays and other databases participating in this offer appear in NEWS 10.

```
=> e us 2006566384
NUMBER OF TERMS TO DISPLAY IS OUT OF RANGE
The total number of terms displayed in a single EXPAND command
must be in the range 5-25.
=> e us 2006566384/apps
                 UA99-95363/PRN
            1
            1
                 UA99-98052573/PRN
E3
            0 --> US 2006566384/AP
E4
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                US 2006566384/PRN
E5
           1
                 US00-10056/AP
                 US00-10070/AP
E.6
           1
                US00-10178/AP
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US00-10197/AP

US00-10220/AP 3 US00-102457/AP 1 E10 E11 US00-10268/AP 1 E12 1 US00-103430/AP

1

1

E7

Ε8

E.9

=> e us 2006-566384/apps 2 US2006-566383/AP E2 1 US2006-566383/PRN 1 --> US2006-566384/AP E.3 1 --> US2UU6-566384/AP 0 US2006-566384/PRN 2 US2006-566385/AP 1 US2006-566387/AP 1 US2006-566388/AP 1 US2006-566389/AP 1 US2006-566390/AP 1 US2006-566390/PRN 2 US2006-566392/AP 1 US2006-566393/AP E.4 E5 Ε6 Ε7 Ε8 E9 E10 E11 E12 1 US2006-566393/AP

=> s us 2006-566384/apps 1 US2006-566384/AP 0 US2006-566384/PRN L11 US 2006-566384/APPS (US2006-566384/AP, PRN)

=> sel rn E1 THROUGH E6 ASSIGNED

=> file registry COST IN U.S. DOLLARS

FULL ESTIMATED COST 6.50 5.80 FILE 'REGISTRY' ENTERED AT 14:25:37 ON 18 NOV 2009

SINCE FILE

ENTRY

TOTAL

SESSION

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8 DICTIONARY FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 26, 2009.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

=> s e1-e61 1044645-80-8/BI (1044645-80-8/RN)1 195875-87-7/BI (195875-87-7/RN) 1 50-36-2/BI (50-36-2/RN)1 54-11-5/BI (54-11-5/RN)1 835651-44-0/BI (835651-44-0/RN)1 87-69-4/BI (87-69-4/RN)L2 6 (1044645-80-8/BI OR 195875-87-7/BI OR 50-36-2/BI OR 54-11-5/BI OR 835651-44-0/BI OR 87-69-4/BI)

=> d hitstr

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

- Index Name, MF, and structure - no RN - All substance data, except sequence data

- FIDE, but only 50 names SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

- Protein sequence data, includes RN SOD

SQD3 - Same as SQD, but 3-letter amino acid codes are used - Protein sequence name information, includes RN SON

EPROP - Table of experimental properties PPROP - Table of predicted properties - EPROP, ETAG, PPROP and SPEC PROP

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations

SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields.

HELP FORMATS -- To see detailed descriptions of the predefined formats.

ENTER DISPLAY FORMAT (IDE):ide

THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N:y

L2 ANSWER 1 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN

RN 1044645-80-8 REGISTRY

ED Entered STN: 29 Aug 2008

CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R,2R,3S,5S)-rel- (CA INDEX NAME)

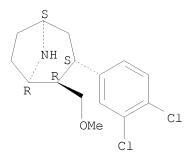
FS STEREOSEARCH

MF C15 H19 Cl2 N O

SR CA

LC STN Files: CA, CAPLUS, TOXCENTER

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)

1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 2

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used SQN - Protein sequence name information, includes RN

EPROP - Table of experimental properties
PPROP - Table of predicted properties
PROP - EPROP, ETAG, PPROP and SPEC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

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IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

OIBIB ---- OBIB, indented with text labels

SBIB ----- BIB, no citations SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats. ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS

THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N:y

L2 ANSWER 2 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN RN 835651-44-0 REGISTRY

ED Entered STN: 22 Feb 2005

CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R,2R,3S,5S)-, (2R,3R)-2,3-dihydroxybutanedioate (1:1) (CA INDEX NAME)

FS STEREOSEARCH

MF C15 H19 C12 N O . C4 H6 O6

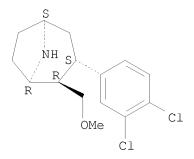
SR CA

LC STN Files: CA, CAPLUS, EMBASE, TOXCENTER, USPATFULL

CM 1

CRN 195875-87-7 CMF C15 H19 C12 N O

Absolute stereochemistry.



CM 2

CRN 87-69-4 CMF C4 H6 O6

Absolute stereochemistry.

2 REFERENCES IN FILE CA (1907 TO DATE)

2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 3

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

```
- Protein sequence data, includes RN
SOD
SQD3
      - Same as SQD, but 3-letter amino acid codes are used
SQN
      - Protein sequence name information, includes RN
      - Table of experimental properties
EPROP
      - Table of predicted properties
PROP
       - EPROP, ETAG, PPROP and SPEC
Any CA File format may be combined with any substance format to
obtain CA references citing the substance. The substance formats
must be cited first. The CA File predefined formats are:
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CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
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IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented
OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) / N: y
    ANSWER 3 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN
L2
     195875-87-7 REGISTRY
RN
ED
     Entered STN: 23 Oct 1997
     8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,
     (1R, 2R, 3S, 5S) - (CA INDEX NAME)
OTHER CA INDEX NAMES:
     8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,
     [1R-(2-endo, 3-exo)]-
FS
     STEREOSEARCH
MF
    C15 H19 C12 N O
CI
    COM
SR
                 CA, CAPLUS, CASREACT, EMBASE, TOXCENTER, USPAT2, USPATFULL
LC
     STN Files:
```

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 6 REFERENCES IN FILE CA (1907 TO DATE) 6 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- => d hitstr 4 'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

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IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

EPROP - Table of experimental properties
PPROP - Table of predicted properties
PROP - EPROP, ETAG, PPROP and SPEC

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CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

```
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
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The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) / N: y
L2
     ANSWER 4 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN
     87-69-4 REGISTRY
RN
ED
     Entered STN: 16 Nov 1984
    Butanedioic acid, 2,3-dihydroxy- (2R,3R)- (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN
     Butanedioic acid, 2,3-dihydroxy-[R-(R^*,R^*)]-
CN
     Tartaric acid, L-(+)-(8CI)
OTHER NAMES:
CN
    (+)-(2R,3R)-Tartaric acid
CN
     (+)-(R,R)-Tartaric acid
CN
     (+)-L-Tartaric acid
CN
     (+)-Tartaric acid
CN
     (2R, 3R) - (+) - Tartaric acid
CN
     (2R, 3R) -2, 3-Dihydroxysuccinic acid
CN
     (2R, 3R) - Tartaric acid
CN
    (R,R)-(+)-Tartaric acid
CN
    (R,R)-Tartaric acid
CN
     1,2-Dihydroxyethane-1,2-dicarboxylic acid
CN
     2,3-Dihydroxybutanedioic acid
CN
     d-\alpha, \beta-Dihydroxysuccinic acid
CN
     d-Tartaric acid
CN
     Dextrotartaric acid
CN
     Dihydroxysuccinic acid
CN
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CN
     L-(+)-Tartaric acid
CN
     L-Tartaric acid
CN
    Natural tartaric acid
     NSC 62778
CN
CN
     Tartaric acid
CN
     Threaric acid
FS
     STEREOSEARCH
     1039646-76-8, 8014-54-8, 8059-77-6, 1336-18-1
DR
MF
     C4 H6 O6
CI
     COM
                  ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS,
     STN Files:
       BIOTECHNO, CA, CABA, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX,
       CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DRUGU, EMBASE, ENCOMPLIT,
       ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, IFICDB, IFIPAT, IFIUDB, IPA,
       MRCK*, MSDS-OHS, NAPRALERT, PIRA, PROMT, PS, RTECS*, SPECINFO,
       SYNTHLINE, TOXCENTER, TULSA, USAN, USPAT2, USPATFULL, USPATOLD
         (*File contains numerically searchable property data)
                     DSL**, EINECS**, TSCA**
     Other Sources:
```

(**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

25806 REFERENCES IN FILE CA (1907 TO DATE)
2072 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
25890 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 5

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except sequence data

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SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used SQN - Protein sequence name information, includes RN

EPROP - Table of experimental properties PPROP - Table of predicted properties PROP - EPROP, ETAG, PPROP and SPEC

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CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

 ${\tt ISTD} \,\, {\tt --} \,\, {\tt STD} \,\, {\tt format} \, , \,\, {\tt indented} \,\,$

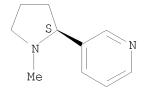
OBIB ----- AN, plus Bibliographic Data (original)

```
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) / N: y
L2
     ANSWER 5 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN
     54-11-5 REGISTRY
RN
ED
     Entered STN: 16 Nov 1984
    Pyridine, 3-[(2S)-1-methyl-2-pyrrolidinyl]- (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN
    Nicotine (8CI)
CN
    Pyridine, 3-(1-methyl-2-pyrrolidinyl)-, (S)-
OTHER NAMES:
CN
    (-)-(S)-Nicotine
CN
     (-)-\beta-Pyridyl-\alpha-N-methylpyrrolidine
CN
     (-)-3-(1-Methyl-2-pyrrolidyl)pyridine
CN
     (-)-Nicotine
CN
     (S)-(-)-Nicotine
CN
     (S)-3-(1-Methyl-2-pyrrolidinyl)pyridine
CN
     (S)-Nicotine
CN
     3-[(2S)-1-Methyl-2-pyrrolidinyl]pyridine
CN
    Exodus
CN
    Flux Maag
CN
    Habitrol
CN
    L-Nicotine
CN
    1-Nicotine
CN
    Nicabate
CN
    Nicoderm
CN
    Nicolan
CN
    Niconil
CN
    Nicopatch
CN
    Nicorette
CN
    Nicotell TTS
CN
    Nicotin
    Nicotinell
CN
CN
    Nicotrol
CN
    Nikofrenon
CN
    Niquitin
CN
    NSC 5065
CN
     Tabazur
     XL All Insecticide
CN
FS
     STEREOSEARCH
DR
     13890-81-8, 13890-82-9, 6912-85-2, 551-13-3, 16760-37-5
MF
    C10 H14 N2
CI
     COM
LC
                  ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*,
     STN Files:
       BIOSIS, BIOTECHNO, CA, CABA, CAPLUS, CASREACT, CBNB, CHEMCATS,
```

CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSCOSEARCH, IMSDRUGNEWS, IMSPATENTS, IMSRESEARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, PHAR, PIRA, PROMT, RTECS*, SPECINFO, TOXCENTER, ULIDAT, USAN, USPAT2, USPATFULL, USPATOLD, VETU

(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
 (**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

22661 REFERENCES IN FILE CA (1907 TO DATE)
395 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
22714 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 6

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except sequence data

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SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used SQN - Protein sequence name information, includes RN

EPROP - Table of experimental properties PPROP - Table of predicted properties PROP - EPROP, ETAG, PPROP and SPEC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

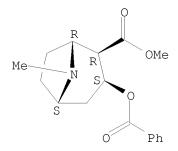
IPC -- International Patent Classification

PATS -- PI, SO

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STD -- BIB, IPC, and NCL
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ISTD -- STD format, indented
OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N:y
L2
     ANSWER 6 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN
RN
     50-36-2 REGISTRY
     Entered STN: 16 Nov 1984
ED
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         (*File contains numerically searchable property data)
                      EINECS**
     Other Sources:
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(**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

15922 REFERENCES IN FILE CA (1907 TO DATE)
333 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
15948 REFERENCES IN FILE CAPLUS (1907 TO DATE)

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(FILE 'HOME' ENTERED AT 14:23:13 ON 18 NOV 2009)

FILE 'REGISTRY' ENTERED AT 14:23:25 ON 18 NOV 2009

FILE 'HCAPLUS' ENTERED AT 14:24:18 ON 18 NOV 2009

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L1 1 S US 2006-566384/APPS SEL RN

FILE 'REGISTRY' ENTERED AT 14:25:37 ON 18 NOV 2009

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L3 1 835651-44-0/RN

=> file hcaplus

COST IN U.S. DOLLARS

SINCE FILE TOTAL
ENTRY SESSION
22.38 28.88

FULL ESTIMATED COST

FILE 'HCAPLUS' ENTERED AT 14:37:58 ON 18 NOV 2009 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2009 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 18 Nov 2009 VOL 151 ISS 21
FILE LAST UPDATED: 17 Nov 2009 (20091117/ED)
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Aug 2009
USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Aug 2009

HCAplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2009.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

During November, try the new LSUS format of legal status information in the CA/CAplus family databases for free! Complete details on the number of free displays and other databases participating in this offer appear in NEWS 10.

=> s 13 L4 2 L3

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THE ESTIMATED COST FOR THIS REQUEST IS 7.78 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N:y

L4 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2008:1070913 HCAPLUS

DOCUMENT NUMBER: 149:315512

TITLE: Hydrochloride salt of an azabicyclo[3.2.1]octane

derivative

INVENTOR(S): Grinter, Trevor John; Moldt, Peter; Watjen, Frank

PATENT ASSIGNEE(S): Glaxo Group Limited, UK; Neurosearch A/S

SOURCE: PCT Int. Appl., 50pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

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IT 835651-44-0

RL: RCT (Reactant); RACT (Reactant or reagent)

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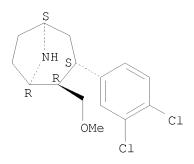
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CRN 195875-87-7 CMF C15 H19 C12 N O

Absolute stereochemistry.



CM 2

CRN 87-69-4 CMF C4 H6 O6

Absolute stereochemistry.

REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2005:120742 HCAPLUS

DOCUMENT NUMBER: 142:183325
TITLE: Preparation of

2-methoxymethyl-3-(3,4-dichlorophenyl)-8-azabicyclo[3.2.1]octane tartrate salts

INVENTOR(S): Frostrup, Brian; Waetjen, Frank; Jensen, Klaus Snej

PATENT ASSIGNEE(S): Neurosearch A/S, Den. SOURCE: PCT Int. Appl., 19 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

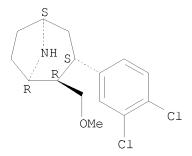
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     BIOL (Biological study); PREP (Preparation); USES (Uses)
        (preparation of methoxymethyldichlorophenyl azabicyclooctane tartrate salts)
RN
     835651-44-0 HCAPLUS
CN
     8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,
     (1R, 2R, 3S, 5S)-, (2R, 3R)-2,3-dihydroxybutanedioate (1:1) (CA INDEX NAME)
     CM
     CRN
          195875-87-7
          C15 H19 C12 N O
     CMF
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Absolute stereochemistry.



CM 2

CRN 87-69-4 CMF C4 H6 O6

Absolute stereochemistry.

OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD

(4 CITINGS)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> file registry
COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
10.63 39.51

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STRUCTURE FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8 DICTIONARY FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8

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http://www.cas.org/support/stngen/stndoc/properties.html

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L5 2 1044645-80-8/RN OR 195875-87-7/RN

=> file hcaplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 0.48 39.99

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FILE COVERS 1907 - 18 Nov 2009 VOL 151 ISS 21
FILE LAST UPDATED: 17 Nov 2009 (20091117/ED)
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Aug 2009
USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Aug 2009

HCAplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2009.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

During November, try the new LSUS format of legal status information in the ${\rm CA/CAplus}$ family databases for free! Complete details on the number of free displays and other databases participating in this offer appear in NEWS 10.

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L7 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2005:120742 HCAPLUS

DOCUMENT NUMBER: 142:183325
TITLE: Preparation of

2-methoxymethyl-3-(3,4-dichlorophenyl)-8-azabicyclo[3.2.1]octane tartrate salts

INVENTOR(S): Frostrup, Brian; Waetjen, Frank; Jensen, Klaus Snej

PATENT ASSIGNEE(S): Neurosearch A/S, Den. SOURCE: PCT Int. Appl., 19 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE

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L12 ANSWER 1 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2008:1070913 HCAPLUS

DOCUMENT NUMBER: 149:315512

TITLE: Hydrochloride salt of an azabicyclo[3.2.1]octane

derivative

INVENTOR(S): Grinter, Trevor John; Moldt, Peter; Watjen, Frank

PATENT ASSIGNEE(S): Glaxo Group Limited, UK; Neurosearch A/S

SOURCE: PCT Int. Appl., 50pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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OTHER SOURCE(S):

CASREACT 149:315512

IT 195875-87-7P

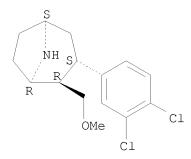
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(hydrochloride salt of an azabicyclo[3.2.1]octane derivative)

RN 195875-87-7 HCAPLUS

CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R,2R,3S,5S)- (CA INDEX NAME)

Absolute stereochemistry.



REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 2 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2007:478903 HCAPLUS

DOCUMENT NUMBER: 146:468550

TITLE: Treatment of diabetes with tropanes

INVENTOR(S): Dugi, Klaus; Berger, Frank; Raschig, Andreas; Reess,

Juergen; Salin, Laurence

PATENT ASSIGNEE(S): Boehringer Ingelheim Pharma G.m.b.H. & Co. K.-G.,

Germany

SOURCE: Eur. Pat. Appl., 19pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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EP	1779	 851						20070502		EP 2005-23794					2	0051	031	
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CA	2627	912			A1		2007	0510		CA 2	2006-	2627	912		2	0061	031	
WO	2007	0515	94		A1		2007	0510		WO 2	2006-	EP10	476		2	0061	031	
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OTHER SOURCE(S): MARPAT 146:468550

IT 195875-87-7

RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)

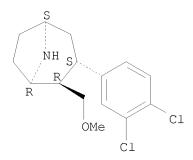
(treatment of diabetes with tropanes)

RN 195875-87-7 HCAPLUS

CN 8-Azabicyclo[3.2.1] octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,

(1R, 2R, 3S, 5S) - (CA INDEX NAME)

Absolute stereochemistry.



OS.CITING REF COUNT: 2 THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD

(2 CITINGS)

REFERENCE COUNT: 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 3 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:971889 HCAPLUS

DOCUMENT NUMBER: 140:13076

TITLE: Triple monoamine reuptake inhibitors for the treatment

of chronic pain

INVENTOR(S): Scheel-Krueger, Jorgen; Blackburn-Munro, Gordon John

PATENT ASSIGNEE(S): Neurosearch A/S, Den. SOURCE: PCT Int. Appl., 20 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.					KIND		DATE		APPLICATION NO.						DATE		
WO	2003	1014	53		A1	_	2003	1211						20030527			
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		FΙ,	FR,	GB,	GR,	HU,	ΙE,	ΙΤ,	LU,	MC,	NL,	PT,	RO,	SE,	SI,	SK,	TR,
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EP	1513	529			A1		2005	0316		EP 2	003-	7248	99		2	0030	527
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CN	1655	787			А		2005	0817		CN 2	003-	8123	76		2	0030	527

JP 2005531594 NZ 536415	T A	20051020 20061027	ΝZ	2004-508810 2003-536415		20030527 20030527
CN 101444509	A	20090603	CN	2008-10176372		20030527
US 20050239824	A1	20051027	US	2004-515275		20041122
US 7459464	В2	20081202				
MX 2004011859	A	20050331	MX	2004-11859		20041129
PRIORITY APPLN. INFO.:			DK	2002-832	A	20020530
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OTHER SOURCE(S): MARPAT 140:13076

IT 195875-87-7

RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL

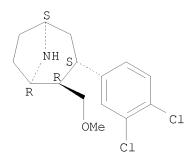
(Biological study); USES (Uses)

(triple monoamine reuptake inhibitors for treatment of chronic pain)

RN 195875-87-7 HCAPLUS

CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R,2R,3S,5S)- (CA INDEX NAME)

Absolute stereochemistry.



OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD

(1 CITINGS)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 4 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:434366 HCAPLUS

DOCUMENT NUMBER: 139:963

TITLE: Tropane derivatives having dopamine reuptake inhibitor

activity for the treatment of ischemic diseases Scheel-Kruger, Jorgen; Ronn, Lars Christian B.

PATENT ASSIGNEE(S): Neurosearch A/S, Den. SOURCE: PCT Int. Appl., 42 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

INVENTOR(S):

PATENT		KIND		DATE		APPLICATION NO.						DATE				
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WO 2003045388				A1 20030605			WO 2002-DK796						20021128			
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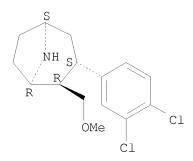
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treatment of ischemic diseases)

195875-87-7 HCAPLUS RN

CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R, 2R, 3S, 5S) - (CA INDEX NAME)

Absolute stereochemistry.



OS.CITING REF COUNT: THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD

(5 CITINGS)

REFERENCE COUNT: 10 THERE ARE 10 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 5 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1997:579720 HCAPLUS

DOCUMENT NUMBER: 127:262895

ORIGINAL REFERENCE NO.: 127:51357a,51360a

TITLE: Preparation of tropane derivatives as inhibitors of

monoamine neurotransmitter re-uptake in the central

nervous system.

Scheel-Kruger, Jorgen; Moldt, Peter; Watjen, Frank INVENTOR(S):

PATENT ASSIGNEE(S): Neurosearch A/S, Den.; Scheel-Kruger, Jorgen; Moldt,

Peter; Watjen, Frank

SOURCE: PCT Int. Appl., 44 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

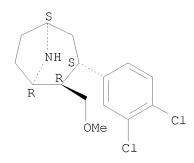
FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9730997	A1	19970828	WO 1997-EP850	19970221

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PRIORITY APPLN. INFO.:
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                         MARPAT 127:262895
OTHER SOURCE(S):
    195875-87-7P
ΙT
     RL: BAC (Biological activity or effector, except adverse); BSU (Biological
     study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
     BIOL (Biological study); PREP (Preparation); USES (Uses)
        (preparation of tropane derivs. as inhibitors of monoamine neurotransmitter
        re-uptake in central nervous system)
RN
     195875-87-7 HCAPLUS
CN
     8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,
     (1R, 2R, 3S, 5S) - (CA INDEX NAME)
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Absolute stereochemistry.



OS.CITING REF COUNT: 27 THERE ARE 27 CAPLUS RECORDS THAT CITE THIS

RECORD (32 CITINGS)

REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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COST IN U.S. DOLLARS SINCE FILE TOTAL

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